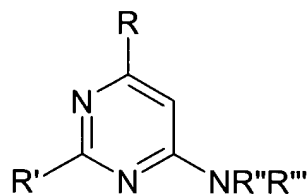


**Amendments to the Claims:**

This listing of claims will replace all prior versions, and listings, of claims in the application:

**Listing of Claims:**

1. (Original) A compound having the structure of general formula (I):



or a salt thereof,

wherein

R represents hydrogen (except when R'=H), (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R' represents hydrogen (except when R=H), (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

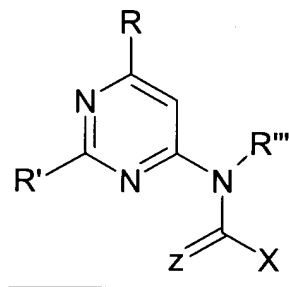
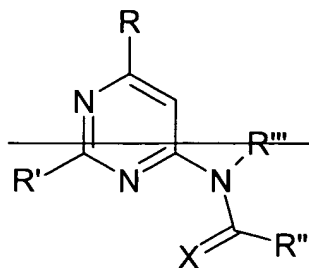
R'' represents hydrogen, acyl, thio-acyl, seleno-acyl, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R''' represents hydrogen, acyl, thio-acyl, seleno-acyl, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R'' and R''' can also together form a substituted or unsubstituted heterocyclic ring or heterocyclic rings;

and n is a number in the range of from 0 to 10.

2. (Currently amended) A compound according to claim 1, having the structure:



or a salt thereof,

wherein

R represents hydrogen (except when R'=H), (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R' represents hydrogen (except when R=H), (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

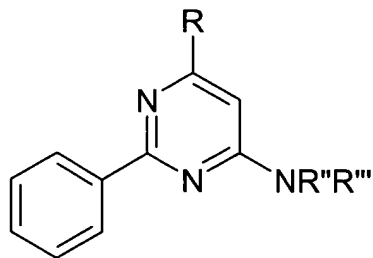
R'' X represents hydrogen, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R''' represents hydrogen, acyl, thio-acyl, seleno-acyl, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

X Z represents oxygen, sulfur or selenium;

and n is a number in the range of from 0 to 10.

3. (Original) A compound according to claim 1, having the structure:



or a salt thereof,

wherein

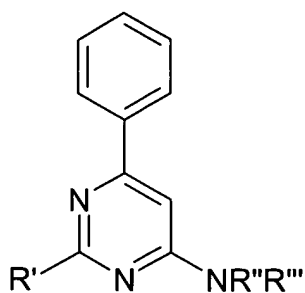
R represents hydrogen, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R'' represents hydrogen, acyl, thio-acyl, seleno-acyl, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R''' represents hydrogen, acyl, thio-acyl, seleno-acyl, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl; and

n is a number in the range of from 0 to 10.

4. (Original) A compound according to claim 1, having the structure:



or a salt thereof,

wherein

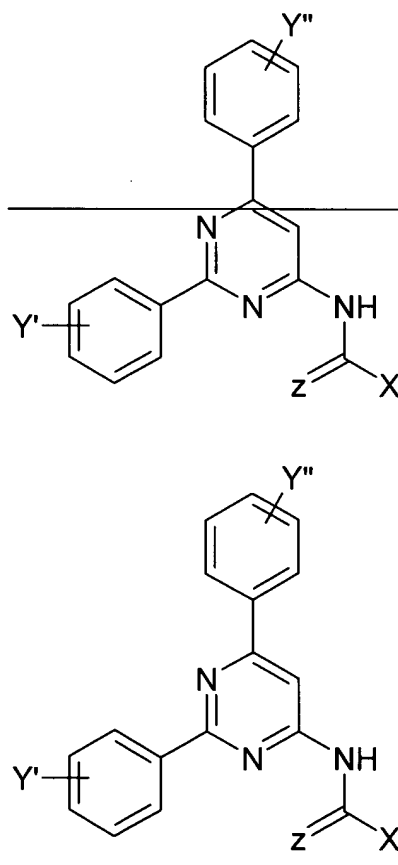
R' represents hydrogen, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R'' represents hydrogen, acyl, thio-acyl, seleno-acyl, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

R''' represents hydrogen, acyl, thio-acyl, seleno-acyl, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl; and

n is a number in the range of from 0 to 10.

5. (Currently amended) The A compound according to ~~any one of claims~~ claim 1, 2 or 4, having the structure:



or a salt thereof,

wherein

R X represents hydrogen, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl;

$R'$   $Y'$  represents hydrogen, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl, alkoxy, thioalkyl, halo,  $NR_1R_2$ ,  $NR_3COR_4$ , or  $NR_5CONR_6R_7$ ;

$R''$   $Y''$  represents hydrogen, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl, alkoxy, thioalkyl, halo,  $NR_1R_2$ ,  $NR_3COR_4$ , or  $NR_5CONR_6R_7$ ; wherein  $R_1$ ,  $R_2$ ,  $R_3$ ,  $R_4$ ,  $R_5$ ,  $R_6$  and  $R_7$  are independently selected from hydrogen, (substituted) alkyl, (substituted) alkenyl, (substituted) alkynyl, or (substituted)  $-(CH_2)_n$ -aryl; and whereby when  $R_1$  and  $R_2$  are in a  $NR_1R_2$  or when  $R_6$  and  $R_7$  are in a  $NR_6R_7$   $R_1$  and  $R_2$  may be linked to form a heterocyclic group;  $X$   $Z$  represents oxygen, sulfur or selenium; and  $n$  is a number in the range of from 0 to 10.

6. (Currently amended) A compound according to claim 1 ~~any one of claims 1-5~~, which compound is selected from the group consisting of N-(2,6-diphenyl-pyrimidin-4-yl)-benzamide, N-(2,6-diphenyl-pyrimidin-4-yl)-4-methoxy-benzamide, N-(2,6-diphenyl-pyrimidin-4-yl)-formamide, N-(2,6-diphenyl-pyrimidin-4-yl)-acetamide, N-(2,6-diphenyl-pyrimidin-4-yl)-propionamide, N-(2,6-diphenyl-pyrimidin-4-yl)-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-isobutyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-3-methyl-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-2-ethyl-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-2-methyl-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-2,2-dimethyl-propionamide, N-(2,6-diphenyl-pyrimidin-4-yl)-3,3-dimethyl-butyramide, cyclopropanecarboxylic acid (2,6-diphenyl-pyrimidin-4-yl)-amide, cyclobutanecarboxylic acid (2,6-diphenyl-pyrimidin-4-yl)-amide, cyclopentanecarboxylic acid (2,6-diphenyl-pyrimidin-4-yl)-amide, cyclohexanecarboxylic acid (2,6-diphenyl-pyrimidin-4-yl)-amide or a salt thereof.

7. (Original) A compound according to claim 6, wherein the compound is selected from the group consisting of N-(2,6-diphenyl-pyrimidin-4-yl)-propionamide, N-(2,6-diphenyl-pyrimidin-4-yl)-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-isobutyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-3-methyl-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-2-ethyl-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-2-methyl-butyramide, N-(2,6-diphenyl-pyrimidin-4-yl)-2,2-dimethyl-propionamide, N-(2,6-diphenyl-pyrimidin-4-yl)-3,3-dimethyl-butyramide,

cyclopentanecarboxylic acid (2,6-diphenyl-pyrimidin-4-yl)-amide, cyclohexanecarboxylic acid (2,6-diphenyl-pyrimidin-4-yl)-amide or a salt thereof.

8. (Currently amended) A compound according to claim 6, ~~which compound comprises~~ wherein the compound is selected from the group consisting of N-(2,6-diphenyl-pyrimidin-4-yl)-2-methyl-butylamide, N-(2,6-diphenyl-pyrimidin-4-yl)-2,2-dimethyl-propionamide, or cyclopentanecarboxylic acid (2,6-diphenyl-pyrimidin-4-yl)-amide or a salt thereof.

9.-10. (Cancelled)

11. (Currently amended) A pharmaceutical composition comprising a pharmaceutically acceptable carrier and a compound according to claim 1 as active ingredient ~~one or more compounds according to any one of claims 1-8.~~

12.-20. (Cancelled)